

AMENDMENTS TO THE SPECIFICATION

Please amend the paragraph at page 3, line 17 as follows

R^5 is selected from a group $NHC(O)OR^9$, $NHC(O)R^9$, $NHS(O)_2R^9$, $C(O)R^9$, $C(O)OR^9$, $S(O)R^9$, $S(O)OR^9$, $S(O)_2OR^9$, $C(O)NR^{10}R^{11}$, $S(O)NR^{10}R^{11}$, $S(O)ONR^{10}R^{11}$ where R^9 , R^{10} or R^{11} are independently selected from hydrogen, optionally ~~substituted~~substituted hydrocarbonyl and optionally substituted heterocyclyl and R^{10} and R^{11} together with the nitrogen atom to which they are attached may additionally form an optionally ~~substituted~~substituted heterocyclic ring which optionally contains further heteroatoms;

Please amend the paragraph at page 4, line 1 as follows

R^6 is hydrogen, optionally ~~substituted~~substituted hydrocarbonyl or optionally substituted heterocyclyl;

Please amend the paragraph at page 4, line 3 as follows

R^7 and R^8 are independently selected from hydrogen, halo, C_{1-4} alkyl, C_{1-4} alkoxy, C_{1-4} alkoxymethyl, di(C_{1-4} alkoxy)methyl, C_{1-4} alkanoyl, trifluoromethyl, cyano, amino, C_{2-5} alkenyl, C_{2-5} alkynyl, a phenyl group, a benzyl group or a 5-6-membered heterocyclic group with 1-3 heteroatoms, selected independently from O, S and N, which heterocyclic group may be aromatic or non-aromatic and may be saturated (linked via a ring carbon or nitrogen atom) or unsaturated (linked via a ring carbon atom), and which phenyl, benzyl or heterocyclic group may bear on one or more ring carbon atoms up to 5 substituents selected from hydroxy, halogeno, C_{1-3} alkyl, C_{1-3} alkoxy, C_{1-3} alkanoyloxy, trifluoromethyl, cyano, amino, nitro, C_{2-4} alkanoyl, C_{1-4} alkanoylamino, C_{1-4} alkoxycarbonyl, C_{1-4} alkylsulphanyl, C_{1-4} alkylsulphinyl, C_{1-4} alkylsulphonyl, carbamoyl, N- C_{1-4} alkylcarbamoyl, N,N-di(C_{1-4} alkyl)carbamoyl, aminosulphonyl, N- C_{1-4} alkylaminosulphonyl, N,N-di(C_{1-4} alkyl)aminosulphonyl, C_{1-4} alkylsulphonylamino, and a saturated heterocyclic group selected from morpholino, thiomorpholino, pyrrolidinyl, piperazinyl, piperidinyl, imidazolidinyl and pyrazolidinyl, which saturated heterocyclic group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno, C_{1-3} alkyl, C_{1-3} alkoxy, C_{1-3} alkanoyloxy, trifluoromethyl, cyano, amino, nitro and C_{1-4} alkoxycarbonyl[[,]]; and

Please amend the paragraph at page 4, line 27 as follows

in the preparation of a medicament for use in the ~~inhibition~~ inhibition of aurora 2 kinase.

Please amend the paragraph at page 5, line 1 as follows

In this specification the term 'alkyl' when used either alone or as a suffix includes straight chained[[,] or branched structures. Unless otherwise stated, these groups may contain up to 10, preferably up to 6 and more preferably up to 4 carbon atoms. Similarly the terms "alkenyl" and "alkynyl" refer to unsaturated straight or branched structures containing for example from 2 to 10, preferably from 2 to 6 carbon atoms. Cyclic moieties such as cycloalkyl, cycloalkenyl and cycloalkynyl are similar in nature but have at least 3 carbon atoms. Terms such as "alkoxy" comprise alkyl groups as is understood in the art.

Please amend the paragraph at page 6, line 1 as follows

The term "functional group" refers to reactive substituents such as nitro, cyano, halo, oxo, $=CR^{78}R^{79}$, $C(O)_xR^{77}$, OR^{77} , $S(O)_yR^{77}$, $NR^{78}R^{79}$, $C(O)NR^{78}R^{79}$, $OC(O)NR^{78}R^{79}$, $=NOR^{77}$, $-NR^{77}C(O)_xR^{78}$, $-NR^{77}CONR^{78}R^{79}$, $-N=CR^{78}R^{79}$, $S(O)_yNR^{78}R^{79}$ or $-NR^{77}S(O)_yR^{78}$ where R^{77} , R^{78} and R^{79} are independently selected from hydrogen, optionally substituted hydrocarbyl, optionally substituted ~~heteroaryl~~ heterocyclyl, or optionally substituted alkoxy, or R^{78} and R^{79} together form an optionally substituted ring which optionally contains further heteroatoms such as oxygen, nitrogen, S, S(O) or S(O)₂, where x is an integer of 1 or 2, y is 0 or an integer of 1-3.

Please amend the paragraph at page 6, line 16 as follows

In particular, optional substituents for hydrocarbyl, ~~heteroaryl~~ heterocyclyl, or alkoxy groups R^{77} , R^{78} and R^{79} include halo, perhaloalkyl, such as trifluoromethyl, mercapto, hydroxy, carboxy, alkoxy, heteroaryl, heteroaryloxy, alkenyloxy, alkynyloxy, alkoxyalkoxy, aryloxy (where the aryl group may be substituted by halo, nitro, or hydroxy), cyano, nitro, amino, mono- or di-alkyl amino, oximino or $S(O)_yR^{90}$ where y is as defined above and R^{90} is a hydrocarbyl group such as alkyl.

Please amend the paragraph at page 7, line 3 as follows

2) $-R^3X^2C(O)R^{19}$ (wherein X^2 represents -O- or $-NR^{20}$ - (in which R^{20} represents hydrogen, or alkyl optionally substituted with a functional group) and R^{19} represents C₁₋₃alkyl, $-NR^{21}R^{22}$ or

-OR²³ (wherein R²¹, R²² and R²³ which may be the same or different each represents hydrogen, or alkyl optionally substituted with a functional group));

Please amend the paragraph at page 7, line 8 as follows

3) -R^bX³R²⁴ (wherein X³ represents -O-, -C(O)-, -S-, -SO-, -SO₂-, -OC(O)-, -NR²⁵C(O)-, -C(O)NR²⁶-, -SO₂NR²⁷-, -NR²⁸SO₂- or -NR²⁹- (wherein R²⁵, R²⁶, R²⁷, R²⁸ and R²⁹ each independently represents hydrogen, or alkyl optionally substituted with a functional group and s is 1 or 2) and R²⁴ represents hydrogen, hydrocarbyl (as defined herein) or a saturated heterocyclic group, wherein the hydrocarbyl or heterocyclic groups may be optionally substituted by one or more functional groups and the heterocyclic groups may additionally be substituted by a hydrocarbyl group);

Please amend the paragraph at page 7, line 27 as follows

9) R³⁷ [[[]] wherein R³⁷ represents a pyridone group, an aryl group or an aromatic heterocyclic group (linked via carbon or nitrogen) with 1-3 heteroatoms selected from O, N and S, which pyridone, aryl or aromatic heterocyclic group may be substituted by one or more functional groups or by a hydrocarbyl group optionally substituted by one or more functional groups or heterocyclyl groups, or by a heterocyclyl group optionally substituted by one or more functional groups or hydrocarbyl groups;

Please amend the paragraph at page 8, line 14 as follows

15) -R^mX⁸R³⁷ (wherein X⁸ represents -O-, -C(O)-, -S-, -SO-, -SO₂-, -OC(O)-, -NR⁵²C(O)-, -C(O)NR⁵³-, -SO₂NR⁵⁴-, -NR⁵⁵SO₂- or -NR⁵⁶- (wherein R⁵², R⁵³, R⁵⁴, R⁵⁵ and R⁵⁶ each independently represents hydrogen, ~~hydrogen~~, or alkyl optionally substituted with a functional group) and R³⁷ is as defined hereinbefore);

Please amend the paragraph at page 8, line 18 as follows

16) -RⁿX⁹R^{n'}R³⁷ (wherein X⁹ represents -O-, -C(O)-, -S-, -SO-, -SO₂-, -OC(O)-, -NR⁵⁷C(O)-, -C(O)NR⁵⁸-, -SO₂NR⁵⁹-, -NR⁶⁰SO₂- or -NR⁶¹- (wherein R⁵⁷, R⁵⁸, R⁵⁹, R⁶⁰ and R⁶¹ each independently represents hydrogen, ~~hydrogen~~, or alkyl optionally substituted with a functional group) and R³⁷ is as defined hereinbefore);

Please amend the paragraph at page 8, line 29 as follows

22) - $R^v R^{62}(R^v)_q(X^9)_r R^{63}$ (wherein X^9 is as defined hereinbefore, q is 0 or 1, r is 0 or 1, and R^{62} is a C_{1-3} alkylene group or a cyclic group selected from divalent cycloalkyl or heterocyclic group, which C_{1-3} alkylene group may be substituted by one or more functional groups and which cyclic group may be substituted by one or more functional groups or by a hydrocarbyl group optionally substituted by one or more functional groups or heterocyclyl groups, or by a heterocyclyl group optionally ~~substituted~~-substituted by one or more functional groups or hydrocarbyl groups; and R^{63} is hydrogen, C_{1-3} alkyl, or a cyclic group selected from cycloalkyl or heterocyclic group, which C_{1-3} alkyl group may be substituted by one or more functional groups and which cyclic group may be substituted by one or more may be substituted by one or more functional groups or by a hydrocarbyl group optionally substituted by one or more functional groups or heterocyclyl groups, or by a heterocyclyl group optionally substituted by one or more functional groups or hydrocarbyl groups);

Please amend the paragraph at page 9, line 11 as follows

and wherein R^a , R^b , $[[]]$, R^c , R^d , R^e , R^f , R^g , R^h , R^i , R^j , R^k , R^l , R^m , R^n , R^o , R^p , R^q , R^r , R^s , R^t , R^u , R^v and R^w are independently selected from C_{1-8} alkylene groups optionally ~~substituted~~-substituted by one or more substituents functional groups $[[,]]$;

Please amend the paragraph at page 12, line 10 as follows

18) C_{2-5} alkenyl which may be unsubstituted or which may be substituted with one or more groups selected from hydroxy, fluoro, amino, C_{1-4} alkylamino, carboxy (and particularly alkyl esters thereof $[[,]]$), N,N-di(C_{1-4} alkyl)amino, aminosulphonyl, N- C_{1-4} alkylaminosulphonyl and N,N-di(C_{1-4} alkyl)aminosulphonyl;

Please amend the paragraph at page 12, line 20 as follows

22) - $R^v R^{62}(R^v)_q(X^9)_r R^{63}$ (wherein X^9 is as defined hereinbefore, q is 0 or 1, r is 0 or 1, and R^{62} is a C_{1-3} alkylene group or a cyclic group selected from cyclopropylene, cyclobutylene, cyclopentylene, cyclohexylene or a 5-6-membered saturated heterocyclic group with 1-2 heteroatoms, selected independently from O, S and N, which C_{1-3} alkylene group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno and C_{1-4} alkoxy and which cyclic group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno, cyano, C_{1-4} cyanoalkyl, C_{1-4} alkyl,

C₁₋₄hydroxyalkyl, C₁₋₄alkoxy, C₁₋₄alkoxyC₁₋₄alkyl, C₁₋₄alkylsulphonylC₁₋₄alkyl, C₁₋₄alkoxycarbonyl, C₁₋₄aminoalkyl, C₁₋₄alkylamino, di(C₁₋₄alkyl)amino, C₁₋₄alkylaminoC₁₋₄alkyl, di(C₁₋₄alkyl)aminoC₁₋₄alkyl, C₁₋₄alkylaminoC₁₋₄alkoxy, di(C₁₋₄alkyl)aminoC₁₋₄alkoxy and a group $-(O-)_f(C_{1-4}alkyl)_g ringD$ (wherein f is 0 or 1, g is 0 or 1 and ring D is a cyclic group selected from C₃₋₆cycloalkyl, aryl or 5-6-membered saturated or unsaturated heterocyclic group with 1-2 heteroatoms, selected independently from O, S and N, which cyclic group may bear one or more substituents selected from halo and C₁₋₄alkyl); and R⁶³ is hydrogen, C₁₋₃alkyl, or a cyclic group selected from cyclopropyl, cyclobutyl, cyclopentyl, cyclohexyl and a 5-6-membered saturated or unsaturated heterocyclic group with 1-2 heteroatoms, selected independently from O, S and N, which C₁₋₃alkyl group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno, C₁₋₄alkoxy and which cyclic group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno, cyano, C₁₋₄cyanoalkyl, C₁₋₄alkyl, C₁₋₄hydroxyalkyl, C₁₋₄alkoxy, C₁₋₄alkoxyC₁₋₄alkyl, C₁₋₄alkylsulphonylC₁₋₄alkyl, C₁₋₄alkoxycarbonyl, C₁₋₄aminoalkyl, C₁₋₄alkylamino, di(C₁₋₄alkyl)amino, C₁₋₄alkylaminoC₁₋₄alkyl, di(C₁₋₄alkyl)aminoC₁₋₄alkyl, C₁₋₄alkylaminoC₁₋₄alkoxy, di(C₁₋₄alkyl)aminoC₁₋₄alkoxy and a group $-(O-)_f(C_{1-4}alkyl)_g ringD$ (wherein f is 0 or 1, g is 0 or 1 and ring D is a cyclic group selected from C₃₋₆cycloalkyl, aryl or 5-6-membered saturated or unsaturated heterocyclic group with 1-2 heteroatoms, selected independently from O, S and N, which cyclic group may bear one or more substituents selected from halo and C₁₋₄alkyl));

Please amend the paragraph at page 13, line 17 as follows

and wherein R^a, R^b, R^{b'}, R^c, R^{c'}, R^d, R^g, R^j, Rⁿ, R^{n'}, R^p, R^{p'}, R^{p'}, R^t, R^u, R^v and R^{v'} are independently selected from C₁₋₈alkylene groups optionally substituted by one or more substituents selected from hydroxy, halogeno, and amino[[,]];

Please amend the paragraph at page 13, line 25 as follows

In particular R¹, R², R³, R⁴ are independently selected from, halo, cyano, nitro, trifluoromethyl, C₁₋₃alkyl, -NR¹³R¹⁴ (wherein R¹³ and R¹⁴, which may be the same or different, each represents hydrogen or C₁₋₃alkyl), or -X¹R¹⁵ [[([)] wherein X¹ represents a direct bond, -O-, -CH₂-, -OCO-, carbonyl, -S-, -SO-, -SO₂-, -NR¹⁶CO-, -CONR¹⁶-, -SO₂NR¹⁶-, -NR¹⁷SO₂- or -NR¹⁸- (wherein R¹⁶, R¹⁷ and R¹⁸ each independently represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl), and R¹⁵ is selected from one of the following groups:

Please amend the paragraph at page 14, line 3 as follows

2') C₁₋₅alkylX²COR¹⁹ (wherein X² represents -O- or -NR²⁰ - (in which R²⁰ represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl) and R¹⁹ represents C₁₋₃alkyl, -NR²¹ or -OR²³ (wherein R²¹, R²² and R²³ which may be the same or different each represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl));

Please amend the paragraph at page 16, line 6 as follows

In particular R¹⁵ is selected from a group of formula (1), (3), (6), (10) or (22) above and preferably selected from groups (1) or (10) above. Particular groups R¹⁵ are those in group (1) above, especially alkyl such as methyl or halo ~~substituted~~-substituted alkyl, or those in group (10) above. In one suitable embodiment, at least one of R² or R³ is a group OC₁₋₅alkylR³⁶ and R³⁶ is a heterocyclic ring such as an N-linked morpholine ring such as 3-morpholinopropoxy.

Please amend the paragraph at page 16, line 20 as follows

Particular examples of R⁶ include ~~h~~-H or heterocyclic groups such as ~~n~~N-morpholino. Preferably however, R⁶ is hydrogen.

Please amend the paragraph at page 17, line 1 as follows

~~heterocycyl~~-heterocyclyl optionally substituted with one or more functional, alkyl, alkenyl or alkynyl groups;

Please amend the paragraph at page 17, line 3 as follows

alkyl optionally substituted by a functional group or a cycloalkyl or heterocyclyl group wherein the cycloalkyl or heterocyclyl group may themselves be optionally substituted with one or more functional or alkyl groups;

Please amend the paragraph at page 17, line 18 as follows

Suitable optionally substituted C₃₋₆cycloalkyl groups R⁹, R¹⁰ and R¹¹ include optionally substituted cyclopropyl, cyclobutyl, cyclopentyl or cyclohexyl any of which may be optionally ~~substituted~~-substituted with for example nitro, halo, carboxy, cyano, C₁₋₄alkyl, C₁₋₄alkoxy, C₁₋₄alkylthio, acetoxy, acetamido, hydroxy, aminosulphonyl, C₁₋₄alkylsulphonyl,

trifluoromethyl, aralkyl, aralkyloxy, or aryl wherein aryl rings in the substituents may themselves be substituted with for example halo, nitro or C₁₋₄alkyl.

Please amend the paragraph at page 17, line 25 as follows

Suitable optionally substituted aralkyl groups R⁹, R¹⁰ and R¹¹ include optionally substituted benzyl, phenylethyl or phenylpropyl, wherein the phenyl ring is optionally substituted with for example up to 5 groups selected from nitro, halo, carboxy, cyano, C₁₋₄alkyl, C₁₋₄alkoxy, C₁₋₄alkylthio, acetoxy, acetamido, hydroxy, aminosulphonyl, C₁₋₄alkylsulphonyl, trifluoromethyl, aralkyl, or aralkyloxy wherein aryl rings in the substituents may themselves be substituted with for example halo, carboxy, trifluoromethyl, nitro or C₁₋₄alkyl and in particular nitro, C₁₋₄alkoxy, halo, hydroxy, trifluoromethyl or carboxy.

Please amend the paragraph at page 18, line 3 as follows

Suitable optionally substituted ~~heterocyclyl~~ heterocyclyl groups R⁹, R¹⁰ and R¹¹ include pyridyl, pyrazine, pyrimidinyl, pyrrolidino, furyl, tetrahydrofuryl, oxazolyl, morpholino, thiadiazole, indolyl, quinoliny, isoquinoliny, pyrazolyl, methylenedioxybenzyl, thiophene, benzothiophene, all of which may be optionally substituted with, for example, one or more groups selected from nitro, halo, carboxy, cyano, C₁₋₄alkyl, C₁₋₄alkoxy, C₁₋₄alkylthio, acetoxy, acetamido hydroxy, aminosulphonyl, C₁₋₄alkylsulphonyl, trifluoromethyl, aralkyl, or aralkyloxy wherein aryl rings in the substituents may themselves be substituted with for example halo, carboxy, trifluoromethyl, nitro or C₁₋₄alkyl; and particularly with C₁₋₄alkyl, halo or nitro.

Please amend the paragraph at page 18, line 19 as follows

Suitable optional substituents for alkenyl or alkynyl groups R⁹, R¹⁰ or R¹¹ include nitro, halo, carboxy, cyano, C₁₋₄alkyl, C₁₋₄alkoxy, C₁₋₄alkylthio, acetoxy, acetamido, hydroxy, aminosulphonyl, C₁₋₄alkylsulphonyl, trifluoromethyl, aralkyl, or aralkyloxy wherein aryl rings in the substituents may themselves be substituted with for example halo, carboxy, trifluoromethyl, nitro or C₁₋₄alkyl. In particular such groups are substituted by aryl such as phenyl, where the aryl ring may itself be substituted with for example halo, nitro, carboxy, or trifluoromethyl.

Please amend the paragraph at page 19, line 10 as follows

in the preparation of a medicament for use in the ~~inhibition~~ inhibition of aurora 2 kinase.

Please amend the paragraph at page 20, line 2 as follows

R⁷ and R⁸ are independently selected from hydrogen, halo, C₁₋₄alkyl, C₁₋₄alkoxy, C₁₋₄alkoxymethyl, di(C₁₋₄alkoxy)methyl, C₁₋₄alkanoyl, trifluoromethyl, cyano, amino, C₂₋₅alkenyl, C₂₋₅alkynyl, a phenyl group, a benzyl group or a 5-6-membered heterocyclic group with 1-3 heteroatoms, selected independently from O, S and N, which heterocyclic group may be aromatic or non-aromatic and may be saturated (linked via a ring carbon or nitrogen atom) or unsaturated (linked via a ring carbon atom), and which phenyl, benzyl or heterocyclic group may bear on one or more ring carbon atoms up to 5 substituents selected from hydroxy, halogeno, C₁₋₃alkyl, C₁₋₃alkoxy, C₁₋₃alkanoyloxy, trifluoromethyl, cyano, amino, nitro, C₂₋₄alkanoyl, C₁₋₄alkanoylamino, C₁₋₄alkoxycarbonyl, C₁₋₄alkylsulphanyl, C₁₋₄alkylsulphinyl, C₁₋₄alkylsulphonyl, carbamoyl, N-C₁₋₄alkylcarbamoyl, N,N-di(C₁₋₄alkyl)carbamoyl, aminosulphonyl, N-C₁₋₄alkylaminosulphonyl, N,N-di(C₁₋₄alkyl)aminosulphonyl, C₁₋₄alkylsulphonylamino, and a saturated heterocyclic group selected from morpholino, thiomorpholino, pyrrolidinyl, piperazinyl, piperidinyl, imidazolidinyl and pyrazolidinyl, which saturated heterocyclic group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno, C₁₋₃alkyl, C₁₋₃alkoxy, C₁₋₃alkanoyloxy, trifluoromethyl, cyano, amino, nitro and C₁₋₄alkoxycarbonyl[[,]]; and

Please amend the paragraph at page 20, line 19 as follows

where R¹, R², R³ and R⁴ are independently selected from, halo, cyano, nitro, trifluoromethyl, C₁₋₃alkyl, -NR¹³R¹⁴ (wherein R¹³ and R¹⁴, which may be the same or different, each represents hydrogen or C₁₋₃alkyl), or -X¹R¹⁵ [[(]] wherein X¹ represents a direct bond, -O-, -CH₂-, -OCO-, carbonyl, -S-, -SO-, -SO₂-, -NR¹⁶CO-, -CONR¹⁶-, -SO₂NR¹⁶-, -NR¹⁷SO₂- or -NR¹⁸- (wherein R¹⁶, R¹⁷ and R¹⁸ each independently represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl), and R¹⁵ is selected from one of the following groups:

Please amend the paragraph at page 20, line 28 as follows

2') C₁₋₅alkylX²COR¹⁹ (wherein X² represents -O- or -NR²⁰ - in which R²⁰ represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl) and R¹⁹ represents C₁₋₃alkyl, -NR[[1]]²¹R²² or -OR²³ (wherein R²¹, R²² and R²³ which may be the same or different each represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl));

Please amend the paragraph at page 22, line 18 as follows

in the preparation of a medicament for use in the ~~inhibition~~inhibition of aurora 2 kinase.

Please amend the paragraph at page 22, line 21 as follows

Particular examples of groups R⁶⁴ include groups listed above for R⁹, and in particular are optionally ~~substituted~~substituted C₁₋₆alkyl, optionally substituted C₂₋₆ alkenyl, optionally substituted phenyl, naphthyl or benzyl, optionally substituted heterocyclyl such as pyridyl[[,]]or furanyl[[,]].

Please amend the paragraph at page 22, line 28 as follows

In particular, the substituents for R⁶⁴ include halo, nitro, optionally substituted C₁₋₆ alkoxy, C₁₋₄alkoxymethyl, di(C₁₋₄alkoxy)methyl, C₁₋₄alkanoyl, trifluoromethyl, cyano, amino, C₂₋₅alkenyl, C₂₋₅alkynyl, a phenyl group, a benzyl group or a 5-6-membered heterocyclic group with 1-3 heteroatoms, selected independently from O, S and N, which heterocyclic group may be aromatic or non-aromatic and may be saturated (linked via a ring carbon or nitrogen atom) or unsaturated (linked via a ring carbon atom), and which phenyl, benzyl or heterocyclic group may bear on one or more ring carbon atoms up to 5 substituents selected from hydroxy, halogeno, C₁₋₃alkyl, C₁₋₃alkoxy, C₁₋₃alkanoyloxy, trifluoromethyl, cyano, amino, nitro, C₂₋₄alkanoyl, C₁₋₄alkanoylamino, C₁₋₄alkoxycarbonyl, C₁₋₄alkylsulphanyl, C₁₋₄alkylsulphinyl, C₁₋₄alkylsulphonyl, carbamoyl, N-C₁₋₄alkylcarbamoyl, N,N-di(C₁₋₄alkyl)carbamoyl, aminosulphonyl, N-C₁₋₄alkylaminosulphonyl, N,N-di(C₁₋₄alkyl)aminosulphonyl, C₁₋₄alkylsulphonylamino, and a saturated heterocyclic group selected from morpholino, thiomorpholino, pyrrolidinyl, piperazinyl, piperidinyl, imidazolidinyl and pyrazolidinyl, which saturated heterocyclic group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno, C₁₋₃alkyl, C₁₋₃alkoxy, C₁₋₃alkanoyloxy, trifluoromethyl, cyano, amino, nitro and C₁₋₄alkoxycarbonyl.

Please amend the paragraph at page 28, line 16 as follows

2') C₁₋₅alkylX²COR¹⁹ (wherein X² represents -O- or -NR²⁰ - in which R²⁰ represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl) and R¹⁹ represents C₁₋₃alkyl, -NR[[1]]²¹R²² or -OR²³ (wherein R²¹, R²² and R²³ which may be the same or different each represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl));

Please amend the paragraph at page 56, line 10 as follows

(iii) where R^1 , R^2 [[⁺]], R^3 , R^4 , R^6 , R^7 and R^8 are all hydrogen, X is oxygen, R^6 is 4-methyl-1-piperazinyl and Z is C(O), R^{64} is other methyl.

Please amend the paragraph at page 56, line 19 as follows

R^7 and R^8 are independently selected from hydrogen, halo, C_{1-4} alkyl, C_{1-4} alkoxy, C_{1-4} alkoxymethyl, di(C_{1-4} alkoxy)methyl, C_{1-4} alkanoyl, trifluoromethyl, cyano, amino, C_{2-5} alkenyl, C_{2-5} alkynyl, a phenyl group, a benzyl group or a 5-6-membered heterocyclic group with 1-3 heteroatoms, selected independently from O, S and N, which heterocyclic group may be aromatic or non-aromatic and may be saturated (linked via a ring carbon or nitrogen atom) or unsaturated (linked via a ring carbon atom), and which phenyl, benzyl or heterocyclic group may bear on one or more ring carbon atoms up to 5 substituents selected from hydroxy, halogeno, C_{1-3} alkyl, C_{1-3} alkoxy, C_{1-3} alkanoyloxy, trifluoromethyl, cyano, amino, nitro, C_{2-4} alkanoyl, C_{1-4} alkanoylamino, C_{1-4} alkoxycarbonyl, C_{1-4} alkylsulphanyl, C_{1-4} alkylsulphonyl, C_{1-4} alkylsulphonyl, carbamoyl, N- C_{1-4} alkylcarbamoyl, N,N-di(C_{1-4} alkyl)carbamoyl, aminosulphonyl, N- C_{1-4} alkylaminosulphonyl, N,N-di(C_{1-4} alkyl)aminosulphonyl, C_{1-4} alkylsulphonylamino, and a saturated heterocyclic group selected from morpholino, thiomorpholino, pyrrolidinyl, piperazinyl, piperidinyl imidazolidinyl and pyrazolidinyl, which saturated heterocyclic group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno, C_{1-3} alkyl, C_{1-3} alkoxy, C_{1-3} alkanoyloxy, trifluoromethyl, cyano, amino, nitro and C_{1-4} alkoxycarbonyl[[₁]]; and

Please amend the paragraph at page 57, line 12 as follows

where R^1 , R^2 , R^3 and R^4 are independently selected from, halo, cyano, nitro, trifluoromethyl, C_{1-3} alkyl, $-NR^{13}R^{14}$ (wherein R^{13} and R^{14} , which may be the same or different, each represents hydrogen or C_{1-3} alkyl), or $-X^1R^{15}$ [[₁]] wherein X^1 represents a direct bond, -O-, -CH₂-, -OCO-, carbonyl, -S-, -SO-, -SO₂-, $-NR^{16}CO$ -, $-CONR^{16}$ -, $-SO_2NR^{16}$ -, $-NR^{17}SO_2$ - or $-NR^{18}$ - (wherein R^{16} , R^{17} and R^{18} each independently represents hydrogen, C_{1-3} alkyl or C_{1-3} alkoxy C_{2-3} alkyl), and R^{15} is selected from one of the following groups:

Please amend the paragraph at page 57, line 21 as follows

2') C₁₋₅alkylX²COR¹⁹ (wherein X² represents -O- or -NR²⁰ - in which R²⁰ represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl) and R¹⁹ represents C₁₋₃alkyl, -NR²¹ or -OR²³ (wherein R²¹, R²² and R²³ which may be the same or different each represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl));

Please amend the paragraph at page 59, line 20 as follows

or a ~~salt~~salt, ester, amide or prodrug thereof[.];

Please amend the paragraph at page 61, line 13 as follows

where X, Y, R¹, R⁴, R⁷, R⁸ are as defined in relation to compound (VIC), R⁶⁵ is as defined ~~in~~ in relation to compound (VIC), and R⁶⁸ and R⁶⁹ are equivalent to R² and R³ in relation to compound (VIC), except that at least one of R⁶⁸ or R⁶⁹ is a group of sub-formula X¹R¹⁵ where R¹⁵ is as defined in relation to compound (VIC), provided that when said one of R⁶⁸ or R⁶⁹ is morpholinopropoxy, the other is not a group of sub-formula (18) as defined in claim 18; and further provided that when when said one of R⁶⁸ or R⁶⁹ is methoxyethoxy, the other is not methoxy.

Please amend the paragraph at page 61, line 20 as follows

In another embodiment, the invention provides a compound of formula (VID) which is of similar structure to (VIA) above but in which X, Y, R¹, R⁴, R⁶, R⁷, R⁸ and R⁶⁵ are as defined in relation to formula (VI), R⁶⁸ is halo, cyano, nitro, trifluoromethyl, C₁₋₃alkyl, -NR¹³R¹⁴ [(I)] wherein R¹³ and R¹⁴ are as defined above in relation to formula (I), or a group -X¹R¹⁵ where X¹ and R¹⁵ are as defined in relation to formula (I) and R¹⁵ is particularly a group of sub group (1) or (10), and R⁶⁹ is C₁₋₆alkoxy optionally substituted by fluorine or a group X¹²R⁷¹ in which X¹² is selected from a group defined for X¹ above, and R⁷¹ is a heterocyclic group, and in particular a 5-6-membered aromatic heterocyclic group (linked via nitrogen) with 1-3 heteroatoms selected from O, N and S; provided that at least one of R⁶⁸ and R⁶⁹ is other than ~~unsubstituted~~unsubstituted methoxy.